

Connecting via Winsock to STN

Welcome to STN International! Enter x:x

LOGINID:ssptansci1625

PASSWORD:

TERMINAL (ENTER 1, 2, 3, OR ?):2

\* \* \* \* \* Welcome to STN International \* \* \* \* \*

NEWS 1 Web Page for STN Seminar Schedule - N. America  
NEWS 2 JAN 02 STN pricing information for 2008 now available  
NEWS 3 JAN 16 CAS patent coverage enhanced to include exemplified  
prophetic substances  
NEWS 4 JAN 28 USPATFULL, USPAT2, and USPATOLD enhanced with new  
custom IPC display formats  
NEWS 5 JAN 28 MARPAT searching enhanced  
NEWS 6 JAN 28 USGENE now provides USPTO sequence data within 3 days  
of publication  
NEWS 7 JAN 28 TOXCENTER enhanced with reloaded MEDLINE segment  
NEWS 8 JAN 28 MEDLINE and LMEEDLINE reloaded with enhancements  
NEWS 9 FEB 08 STN Express, Version 8.3, now available  
NEWS 10 FEB 20 PCI now available as a replacement to DPCI  
NEWS 11 FEB 25 IFIREF reloaded with enhancements  
NEWS 12 FEB 25 IMSPRODUCT reloaded with enhancements  
NEWS 13 FEB 29 WPINDEX/WPIDS/WPIX enhanced with ECLA and current  
U.S. National Patent Classification  
NEWS 14 MAR 31 IFICDB, IFIPAT, and IFIUDB enhanced with new custom  
IPC display formats  
NEWS 15 MAR 31 CAS REGISTRY enhanced with additional experimental  
spectra  
NEWS 16 MAR 31 CA/Caplus and CASREACT patent number format for U.S.  
applications updated  
NEWS 17 MAR 31 LPCI now available as a replacement to LDPCI  
NEWS 18 MAR 31 EMBASE, EMBAL, and LEMBASE reloaded with enhancements  
NEWS 19 APR 04 STN AnaVist, Version 1, to be discontinued  
NEWS 20 APR 15 WPIDS, WPINDEX, and WPIX enhanced with new  
predefined hit display formats  
NEWS 21 APR 28 EMBASE Controlled Term thesaurus enhanced  
NEWS 22 APR 28 IMSRESEARCH reloaded with enhancements  
  
NEWS EXPRESS FEBRUARY 08 CURRENT WINDOWS VERSION IS V8.3,  
AND CURRENT DISCOVER FILE IS DATED 20 FEBRUARY 2008  
  
NEWS HOURS STN Operating Hours Plus Help Desk Availability  
NEWS LOGIN Welcome Banner and News Items  
NEWS IPC8 For general information regarding STN implementation of IPC 8

Enter NEWS followed by the item number or name to see news on that  
specific topic.

All use of STN is subject to the provisions of the STN Customer  
agreement. Please note that this agreement limits use to scientific

research. Use for software development or design or implementation of commercial gateways or other similar uses is prohibited and may result in loss of user privileges and other penalties.

\*\*\*\*\* STN Columbus \*\*\*\*\*

FILE 'HOME' ENTERED AT 15:39:46 ON 04 MAY 2008

=> fil reg

COST IN U.S. DOLLARS

SINCE FILE

TOTAL

ENTRY

SESSION

FULL ESTIMATED COST

0.21

0.21

FILE 'REGISTRY' ENTERED AT 15:39:56 ON 04 MAY 2008

USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.

PLEASE SEE "HELP USAGETERMS" FOR DETAILS.

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Property values tagged with IC are from the ZIC/VINITI data file provided by InfoChem.

STRUCTURE FILE UPDATES: 2 MAY 2008 HIGHEST RN 1019057-98-7

DICTIONARY FILE UPDATES: 2 MAY 2008 HIGHEST RN 1019057-98-7

New CAS Information Use Policies, enter HELP USAGETERMS for details.

TSCA INFORMATION NOW CURRENT THROUGH January 9, 2008.

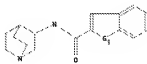
Please note that search-term pricing does apply when conducting SmartSELECT searches.

REGISTRY includes numerically searchable data for experimental and predicted properties as well as tags indicating availability of experimental property data in the original document. For information on property searching in REGISTRY, refer to:

<http://www.cas.org/support/stngen/stdoc/properties.html>

=>

Uploading C:\Program Files\Stnexp\Queries\10522611.str



chain nodes :

9 10 11

ring nodes :

1 2 3 4 5 6 7 8 12 13 14 15 16 17 18 19 20

chain bonds :

4-9 9-10 10-11 10-12

ring bonds :

1-2 1-6 2-3 3-4 3-7 4-5 5-6 6-8 7-8 12-13 12-16 13-14 14-17 14-15 15-16 15-20 17-18 18-19 19-20  
 exact/norm bonds :  
 1-2 1-6 2-3 3-4 3-7 4-5 4-9 5-6 6-8 7-8 9-10 10-11 10-12 12-13 12-16 13-14 15-16  
 normalized bonds :  
 14-17 14-15 15-20 17-18 18-19 19-20

G1:O,S

Match level :

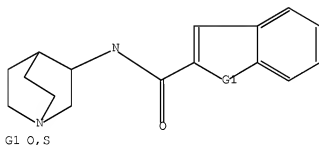
1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:CLASS 10:CLASS  
 11:CLASS 12:Atom 13:Atom 14:Atom 15:Atom 16:Atom 17:Atom 18:Atom 19:Atom  
 20:Atom

L1 STRUCTURE UPLOADED

=> d l1

L1 HAS NO ANSWERS

L1 STR



Structure attributes must be viewed using STN Express query preparation.

=> s sss sam l1

SAMPLE SEARCH INITIATED 15:40:18 FILE 'REGISTRY'

SAMPLE SCREEN SEARCH COMPLETED - 42 TO ITERATE

100.0% PROCESSED 42 ITERATIONS

27 ANSWERS

SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE \*\*COMPLETE\*\*  
 BATCH \*\*COMPLETE\*\*

PROJECTED ITERATIONS: 452 TO 1228

PROJECTED ANSWERS: 229 TO 851

L2 27 SEA SSS SAM L1

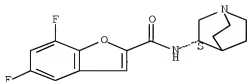
=> d scan

L2 27 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN

IN 2-Benzofurancarboxamide, N-(3S)-1-azabicyclo[2.2.2]oct-3-yl-5,7-difluoro-

MF C16 H16 F2 N2 O2

Absolute stereochemistry.

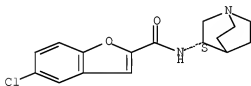


\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):4

L2 27 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN  
IN 2-Benzofurancarboxamide, N-(3S)-1-azabicyclo[2.2.2]oct-3-yl-5-chloro-,  
monohydrochloride (9CI)  
MF C16 H17 Cl N2 O2 . Cl H

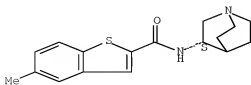
Absolute stereochemistry.



● HCl

```
L2 27 ANSWERS  REGISTRY  COPYRIGHT 2008 ACS on STN
IN Benzo[b]thiophene-2-carboxamide, N-(3S)-1-azabicyclo[2.2.2]oct-3-yl-5-
   methyl-
MF C17 H20 N2 O S
CI COM
```

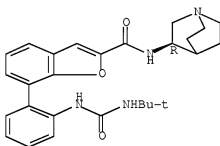
Absolute stereochemistry.



\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

L2 27 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN  
IN 2-Benzofurancarboxamide, N-(3R)-1-azabicyclo[2.2.2]oct-3-yl-7-[2-[[[(1,1-dimethylethyl)amino]carbonyl]amino]phenyl]-  
MF C27 H32 N4 O3  
CI COM

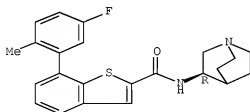
Absolute stereochemistry.



\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

L2 27 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN  
IN Benzo[b]thiophene-2-carboxamide, N-(3R)-1-azabicyclo[2.2.2]oct-3-yl-7-(5-fluoro-2-methylphenyl)-  
MF C23 H23 F N2 O S  
CI COM

Absolute stereochemistry.



\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):  
HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):0

=> s sss full 11

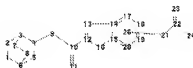
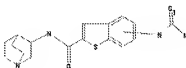
FULL SEARCH INITIATED 15:41:07 FILE 'REGISTRY'  
FULL SCREEN SEARCH COMPLETED - 831 TO ITERATE

100.0% PROCESSED 831 ITERATIONS 647 ANSWERS  
SEARCH TIME: 00.00.01

L3 647 SEA SSS FUL L1

=> save l3 fles10522611/a  
ANSWER SET L3 HAS BEEN SAVED AS 'FLES10522611/A'

=>  
Uploading C:\Program Files\Stnexp\Queries\10522611A.str



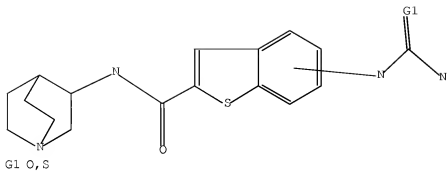
chain nodes :  
9 10 11 23  
ring nodes :  
1 2 3 4 5 6 7 8 12 13 14 15 16 17 18 19 20  
ring/chain nodes :  
21 22 24  
chain bonds :  
4-9 9-10 10-11 10-12 21-22 22-23 22-24  
ring bonds :  
1-2 1-6 2-3 3-4 3-7 4-5 5-6 6-8 7-8 12-13 12-16 13-14 14-17 14-15 15-16 15-20 17-18 18-19 19-20  
exact/norm bonds :  
1-2 1-6 2-3 3-4 3-7 4-5 4-9 5-6 6-8 7-8 9-10 10-11 12-13 12-16 13-14 15-16 21-22 22-23 22-24  
exact bonds :  
10-12  
normalized bonds :  
14-17 14-15 15-20 17-18 18-19 19-20

G1:O,S

Match level :  
1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:CLASS 10:CLASS  
11:CLASS 12:Atom 13:Atom 14:Atom 15:Atom 16:Atom 17:Atom 18:Atom 19:Atom  
20:Atom 21:CLASS 22:CLASS 23:CLASS 24:CLASS 26:Atom

L4 STRUCTURE UPLOADED

=> D L4  
L4 HAS NO ANSWERS  
L4 STR



Structure attributes must be viewed using STN Express query preparation.

```
=> S SSS SAM SUBSET=L3
ENTER LOGIC EXPRESSION, QUERY NAME, OR (END):END
SEARCH ENDED BY USER
```

```
=> S SSS SAM SUBSET=L3 L4
SAMPLE SUBSET SEARCH INITIATED 15:47:10 FILE 'REGISTRY'
SAMPLE SUBSET SCREEN SEARCH COMPLETED - 2 TO ITERATE
```

```
100.0% PROCESSED 2 ITERATIONS 2 ANSWERS
SEARCH TIME: 00.00.01
```

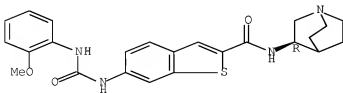
```
PROJECTIONS (WITHIN SPECIFIED SUBSET): ONLINE **COMPLETE**
PROJECTED ITERATIONS (WITHIN SPECIFIED SUBSET): 2 TO 124
PROJECTED ANSWERS (WITHIN SPECIFIED SUBSET): 2 TO 124
```

```
L5 2 SEA SUB=L3 SSS SAM L4
```

```
=> D SCAN
```

```
L5 2 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
IN Benzo[b]thiophene-2-carboxamide, N-(3R)-1-azabicyclo[2.2.2]oct-3-yl-6-
[[[(2-methoxyphenyl)amino]carbonyl]amino]-, monohydrochloride (9CI)
MF C24 H26 N4 O3 S . Cl H
```

Absolute stereochemistry.



\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):0

=> S SSS FULL SUBSET=L3 L4

FULL SUBSET SEARCH INITIATED 15:47:28 FILE 'REGISTRY'

FULL SUBSET SCREEN SEARCH COMPLETED - 44 TO ITERATE

100.0% PROCESSED 44 ITERATIONS

42 ANSWERS

SEARCH TIME: 00.00.01

L6 42 SEA SUB=L3 SSS FUL L4

=> FIL CAPLU

COST IN U.S. DOLLARS

SINCE FILE

TOTAL

ENTRY

SESSION

FULL ESTIMATED COST

225.98

226.19

FILE 'CAPLUS' ENTERED AT 15:47:38 ON 04 MAY 2008

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FILE COVERS 1907 - 4 May 2008 VOL 148 ISS 19

FILE LAST UPDATED: 2 May 2008 (20080502/ED)

Effective October 17, 2005, revised CAS Information Use Policies apply. They are available for your review at:

<http://www.cas.org/infopolicy.html>

=> S L6

L7 1 L6

=> D IBIB ABS HITSTR

L7 ANSWER 1 OF 1 CAPLUS COPYRIGHT 2008 ACS on STN

ACCESSION NUMBER: 2004:117115 CAPLUS Full-text

DOCUMENT NUMBER: 140:181328

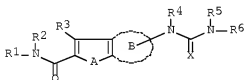
TITLE: Preparation of benzothiophenes, benzofuranes and indoles for improvement of cognition, attention achievement, learning achievement and/or memory achievement

INVENTOR(S): Flessner, Timo; Boess, Frank-Gerhard; Erb, Christina; Hafner, Frank-Thorsten; Schnizler, Katrin; Lang, Dieter; Luithle, Joachim; Van Kampen, Marja; Van der Staay, Franz-Josef



PATENT ASSIGNEE(S): Bayer AG, Germany  
 SOURCE: Ger. Offen., 35 pp.  
 CODEN: GWXXBX  
 DOCUMENT TYPE: Patent  
 LANGUAGE: German  
 FAMILY ACC. NUM. COUNT: 1  
 PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
DE 10234424	A1	20040212	DE 2002-10234424	20020729
CA 2494352	A1	20040212	CA 2003-2494352	20030714
WO 2004013136	A1	20040212	WO 2003-EP7588	20030714
W:	AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NI, NO, NZ, OM, PH, PL, PT, RO, RU, SC, SD, SE, SG, SK, SL, TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VC, VN, YU, ZA, ZM, ZW			
RW:	GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IT, LU, MC, NL, PT, RO, SE, SI, SK, TR, BF, BJ, BR, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG			
AU 2003254343	A1	20040223	AU 2003-254343	20030714
EP 1527073	A1	20050504	EP 2003-766162	20030714
EP 1527073	B1	20070822		
R:	AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, SI, LT, LV, FI, RO, MK, CY, AL, TR, BG, CZ, EE, HU, SK			
JP 2005538111	T	20051215	JP 2004-525191	20030714
ES 2291694	T3	20080301	ES 2003-766162	20030714
US 20060106096	A1	20060518	US 2005-522611	20051215
PRIORITY APPLN. INFO.:			DE 2002-10234424	A 20020729
			WO 2003-EP7588	W 20030714
OTHER SOURCE(S):	MARPAT 140:181328			
GI				



I

AB Title compds. [I; R1 = 1-azabicyclo[2.2.2]oct-3-yl; R2 = H, alkyl; R3 = H, halo, amino, OH, C1-6 alkyl; R4 = H, (substituted) C1-6 alkyl; R5 = H, C1-6 alkyl; or R4R5 = (substituted) 5-6 membered heterocyclyl; R6 = H, alkyl, cycloalkyl, aryl, (substituted) 5-6 membered heterocyclyl, etc.; NR5R6 = (substituted) 3-8 membered heterocyclyl; A = O, N, S; X = O, S; B = (substituted) benzo or pyrido ring], were prepared Thus, 6-amino-N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-1-benzothiophene-2-carboxamide dihydrochloride (preparation given) in THF was reacted with Et3N and 2,6-difluorophenyl isocyanate for 18 h at room temperature to give 75.9% N-[(3R)-1-azabicyclo[2.2.2]oct-3-yl]-6-[(2,6-difluorophenyl)aminocarbonyl]amino]-1-benzothiophene-2-carboxamide hydrochloride. The latter showed affinity to  $\alpha 7$ -NACHR by inhibiting [3H]methyllycaconite binding in rat brain membranes with  $K_i = 0.4$  nM.

IT 657401-24-6P 657401-25-7P 657401-26-8P  
 657401-27-9P 657401-28-0P 657401-29-1P  
 657401-30-4P 657401-31-5P 657401-32-6P  
 657401-32-7P 657401-34-8P 657401-36-0P  
 657401-37-1P 657401-38-2P 657401-39-3P  
 657401-40-6P 657401-41-7P 657401-42-8P  
 657401-43-9P 657401-44-0P 657401-45-1P

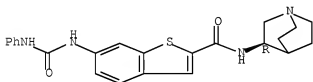
RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU  
 (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES  
 (Uses)

(preparation of benzothiophenes, benzofuranes and indoles for improvement of  
 cognition, attention achievement, learning achievement and/or memory  
 achievement)

RN 657401-24-6 CAPLUS

CN Benzo[b]thiophene-2-carboxamide, N-(3R)-1-azabicyclo[2.2.2]oct-3-yl-6-  
 [[[phenylamino]carbonyl]amino]-, monohydrochloride (9CI) (CA INDEX NAME)

Absolute stereochemistry.

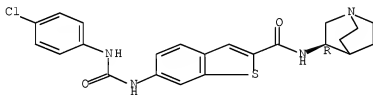


● HCl

RN 657401-25-7 CAPLUS

CN Benzo[b]thiophene-2-carboxamide, N-(3R)-1-azabicyclo[2.2.2]oct-3-yl-6-  
 [[[(4-chlorophenyl)amino]carbonyl]amino]-, monohydrochloride (9CI) (CA  
 INDEX NAME)

Absolute stereochemistry.

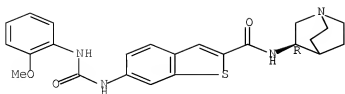


● HCl

RN 657401-26-8 CAPLUS

CN Benzo[b]thiophene-2-carboxamide, N-(3R)-1-azabicyclo[2.2.2]oct-3-yl-6-  
 [[[(2-methoxyphenyl)amino]carbonyl]amino]-, monohydrochloride (9CI) (CA  
 INDEX NAME)

Absolute stereochemistry.

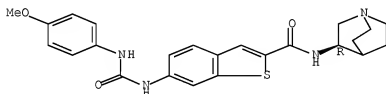


● HCl

RN 657401-27-9 CAPLUS

CN Benzo[b]thiophene-2-carboxamide, N-(3R)-1-azabicyclo[2.2.2]oct-3-yl-6-[[[(4-methoxyphenyl)amino]carbonyl]amino]-, monohydrochloride (9CI) (CA INDEX NAME)

Absolute stereochemistry.

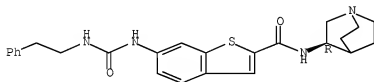


● HCl

RN 657401-28-0 CAPLUS

CN Benzo[b]thiophene-2-carboxamide, N-(3R)-1-azabicyclo[2.2.2]oct-3-yl-6-[[[(2-phenylethyl)amino]carbonyl]amino]-, monohydrochloride (9CI) (CA INDEX NAME)

Absolute stereochemistry.

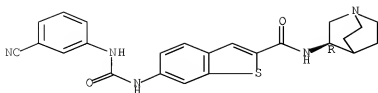


● HCl

RN 657401-29-1 CAPLUS

CN Benzo[b]thiophene-2-carboxamide, N-(3R)-1-azabicyclo[2.2.2]oct-3-yl-6-[[[(3-cyanophenyl)amino]carbonyl]amino]-, monohydrochloride (9CI) (CA INDEX NAME)

Absolute stereochemistry.

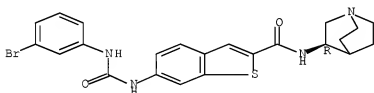


● HCl

RN 657401-30-4 CAPLUS

CN Benzo[b]thiophene-2-carboxamide, N-(3R)-1-azabicyclo[2.2.2]oct-3-yl-6-[[[(3-bromophenyl)amino]carbonyl]amino]-, monohydrochloride (9CI) (CA INDEX NAME)

Absolute stereochemistry.

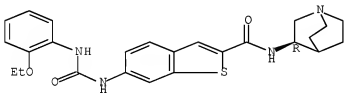


● HCl

RN 657401-31-5 CAPLUS

CN Benzo[b]thiophene-2-carboxamide, N-(3R)-1-azabicyclo[2.2.2]oct-3-yl-6-[[[(2-ethoxyphenyl)amino]carbonyl]amino]-, monohydrochloride (9CI) (CA INDEX NAME)

Absolute stereochemistry.

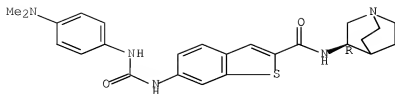


● HCl

RN 657401-32-6 CAPLUS

CN Benzo[b]thiophene-2-carboxamide, N-(3R)-1-azabicyclo[2.2.2]oct-3-yl-6-[[[(4-(dimethylamino)phenyl)amino]carbonyl]amino]-, monohydrochloride (9CI) (CA INDEX NAME)

Absolute stereochemistry.

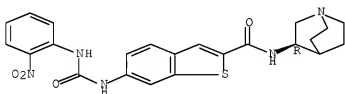


● HCl

RN 657401-33-7 CAPLUS

CN Benzo[b]thiophene-2-carboxamide, N-(3R)-1-azabicyclo[2.2.2]oct-3-yl-6-  
[[[(2-nitrophenyl)amino]carbonyl]amino]-, monohydrochloride (9CI) (CA  
INDEX NAME)

Absolute stereochemistry.

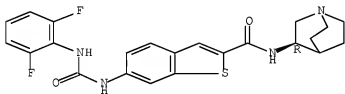


● HCl

RN 657401-34-8 CAPLUS

CN Benzo[b]thiophene-2-carboxamide, N-(3R)-1-azabicyclo[2.2.2]oct-3-yl-6-  
[[[(2,6-difluorophenyl)amino]carbonyl]amino]-, monohydrochloride (9CI)  
(CA INDEX NAME)

Absolute stereochemistry.

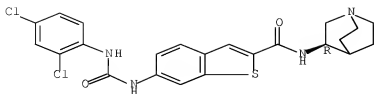


● HCl

RN 657401-36-0 CAPLUS

CN Benzo[b]thiophene-2-carboxamide, N-(3R)-1-azabicyclo[2.2.2]oct-3-yl-6-  
[[[(2,4-dichlorophenyl)amino]carbonyl]amino]-, monohydrochloride (9CI)  
(CA INDEX NAME)

Absolute stereochemistry.

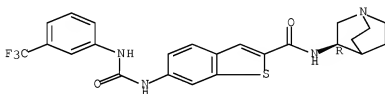


● HCl

RN 657401-37-1 CAPLUS

CN Benzo[b]thiophene-2-carboxamide, N-((3R)-1-azabicyclo[2.2.2]oct-3-yl-6-[[[3-(trifluoromethyl)phenyl]amino]carbonyl]amino)-, monohydrochloride (9CI) (CA INDEX NAME)

Absolute stereochemistry.

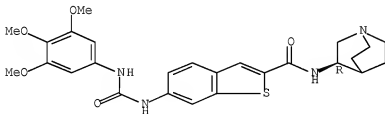


● HCl

RN 657401-38-2 CAPLUS

CN Benzo[b]thiophene-2-carboxamide, N-((3R)-1-azabicyclo[2.2.2]oct-3-yl-6-[[[3,4,5-trimethoxyphenyl]amino]carbonyl]amino)-, monohydrochloride (9CI) (CA INDEX NAME)

Absolute stereochemistry.

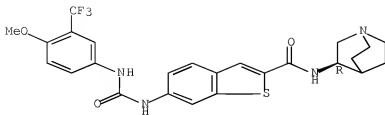


● HCl

RN 657401-39-3 CAPLUS

CN Benzo[b]thiophene-2-carboxamide, N-((3R)-1-azabicyclo[2.2.2]oct-3-yl-6-[[[4-methoxy-3-(trifluoromethyl)phenyl]amino]carbonyl]amino)-, monohydrochloride (9CI) (CA INDEX NAME)

Absolute stereochemistry.

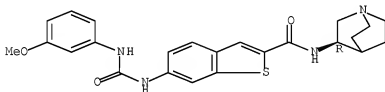


● HCl

RN 657401-40-6 CAPLUS

CN Benzo[b]thiophene-2-carboxamide, N-(3R)-1-azabicyclo[2.2.2]oct-3-yl-6-[[[(3-methoxyphenyl)amino]carbonyl]amino]-, monohydrochloride (9CI) (CA INDEX NAME)

Absolute stereochemistry.

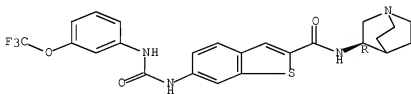


● HCl

RN 657401-41-7 CAPLUS

CN Benzo[b]thiophene-2-carboxamide, N-(3R)-1-azabicyclo[2.2.2]oct-3-yl-6-[[[(3-(trifluoromethoxy)phenyl)amino]carbonyl]amino]-, monohydrochloride (9CI) (CA INDEX NAME)

Absolute stereochemistry.

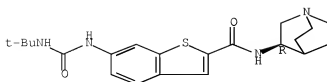


● HCl

RN 657401-42-8 CAPLUS

CN Benzo[b]thiophene-2-carboxamide, N-(3R)-1-azabicyclo[2.2.2]oct-3-yl-6-[[[(1,1-dimethylethyl)amino]carbonyl]amino]-, monohydrochloride (9CI) (CA INDEX NAME)

Absolute stereochemistry.

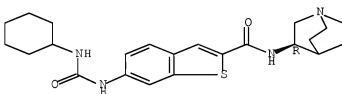


● HCl

RN 657401-43-9 CAPLUS

CN Benzo[b]thiophene-2-carboxamide, N-(3R)-1-azabicyclo[2.2.2]oct-3-yl-6-[[[(cyclohexylamino)carbonyl]amino]-, monohydrochloride (9CI) (CA INDEX NAME)

Absolute stereochemistry.

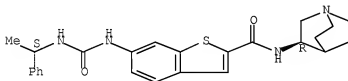


● HCl

RN 657401-44-0 CAPLUS

CN Benzo[b]thiophene-2-carboxamide, N-(3R)-1-azabicyclo[2.2.2]oct-3-yl-6-[[[(1S)-1-phenylethyl]amino]carbonyl]amino]-, monohydrochloride (9CI) (CA INDEX NAME)

Absolute stereochemistry.



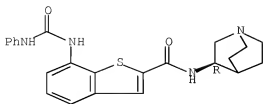
● HCl

RN 657401-45-1 CAPLUS

CN Benzo[b]thiophene-2-carboxamide, N-(3R)-1-azabicyclo[2.2.2]oct-3-yl-7-[[[phenylamino]carbonyl]amino]-, monohydrochloride (9CI) (CA INDEX NAME)

Absolute stereochemistry.





● HCl

=> LOGOFF Y

COST IN U.S. DOLLARS

SINCE FILE

TOTAL

ENTRY

SESSION

FULL ESTIMATED COST

5.93

232.12

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)

SINCE FILE

TOTAL

ENTRY

SESSION

CA SUBSCRIBER PRICE

-0.80

-0.80

STN INTERNATIONAL LOGOFF AT 15:47:57 ON 04 MAY 2008